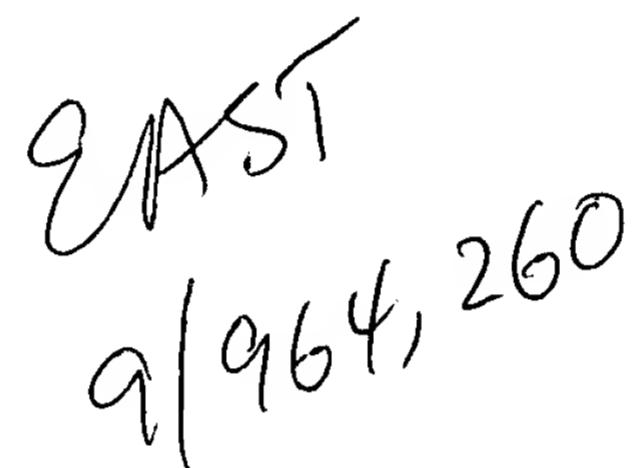


L Number	Hits	Search Text	DB	Time stamp
1	13	pyrimido with (isoquinolin or 'isoquinolin-4-one')	USPAT; US-PGPUB	2003/08/15 14:53
2	29	treqinsin	USPAT; US-PGPUB	2003/08/15 14:53
3	38	(pyrimido with (isoquinolin or 'isoquinolin-4-one')) or treqinsin	USPAT; US-PGPUB	2003/08/15 14:53


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NEWS 14 Apr 21 New current-awareness alert (SDI) frequency in WPIDS/WPINDEX/WPIX
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NEWS 16 May 05 Pharmacokinetic information and systematic chemical names added to PHAR
NEWS 17 May 15 MEDLINE file segment of TOXCENTER reloaded
NEWS 18 May 15 Supporter information for ENCOMPPAT and ENCOMPLIT updated
NEWS 19 May 19 Simultaneous left and right truncation added to WSCA
NEWS 20 May 19 RAPRA enhanced with new search field, simultaneous left and right truncation
NEWS 21 Jun 06 Simultaneous left and right truncation added to CBNB
NEWS 22 Jun 06 PASCAL enhanced with additional data
NEWS 23 Jun 20 2003 edition of the FSTA Thesaurus is now available
NEWS 24 Jun 25 HSDB has been reloaded
NEWS 25 Jul 16 Data from 1960-1976 added to RDISCLOSURE
NEWS 26 Jul 21 Identification of STN records implemented
NEWS 27 Jul 21 Polymer class term count added to REGISTRY
NEWS 28 Jul 22 INPADOC: Basic index (/BI) enhanced; Simultaneous Left and Right Truncation available
NEWS 29 AUG 05 New pricing for EUROPATFULL and PCTFULL effective August 1, 2003
NEWS 30 AUG 13 Field Availability (/FA) field enhanced in BEILSTEIN
NEWS 31 AUG 15 PATDPAFULL: one FREE connect hour, per account, in September 2003
NEWS 32 AUG 15 PCTGEN: one FREE connect hour, per account, in September 2003
NEWS 33 AUG 15 RDISCLOSURE: one FREE connect hour, per account, in September 2003
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DICTIONARY FILE UPDATES: 13 AUG 2003 HIGHEST RN 566135-25-9

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Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

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L1 STRUCTURE UPLOADED

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L1 HAS NO ANSWERS
L1 ST

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Structure attributes must be viewed using STN Express query preparation.

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FULL SEARCH INITIATED 14:20:30 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 14 TO ITERATE
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09/ 964,260

100.0% PROCESSED 14 ITERATIONS
SEARCH TIME: 00.00.01

14 ANSWERS

L2 14 SEA SSS FUL L1

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FULL ESTIMATED COST

SINCE FILE ENTRY	TOTAL SESSION
148.15	148.36

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FILE LAST UPDATED: 14 Aug 2003 (20030814/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 12
L3 1 L2

=> d 13 1- ibib abs hitstr
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L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2003 ACS on STN
ACCESSION NUMBER: 2000:707163 CAPLUS
DOCUMENT NUMBER: 133:266869
TITLE: Preparation of 2-aryliminopyrimido[6,1-a]isoquinolin-4-ones as phosphodiesterase inhibitors.
INVENTOR(S): Oxford, Alexander William; Jack, David
PATENT ASSIGNEE(S): Vanguard Medica Ltd., UK
SOURCE: PCT Int. Appl., 77 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

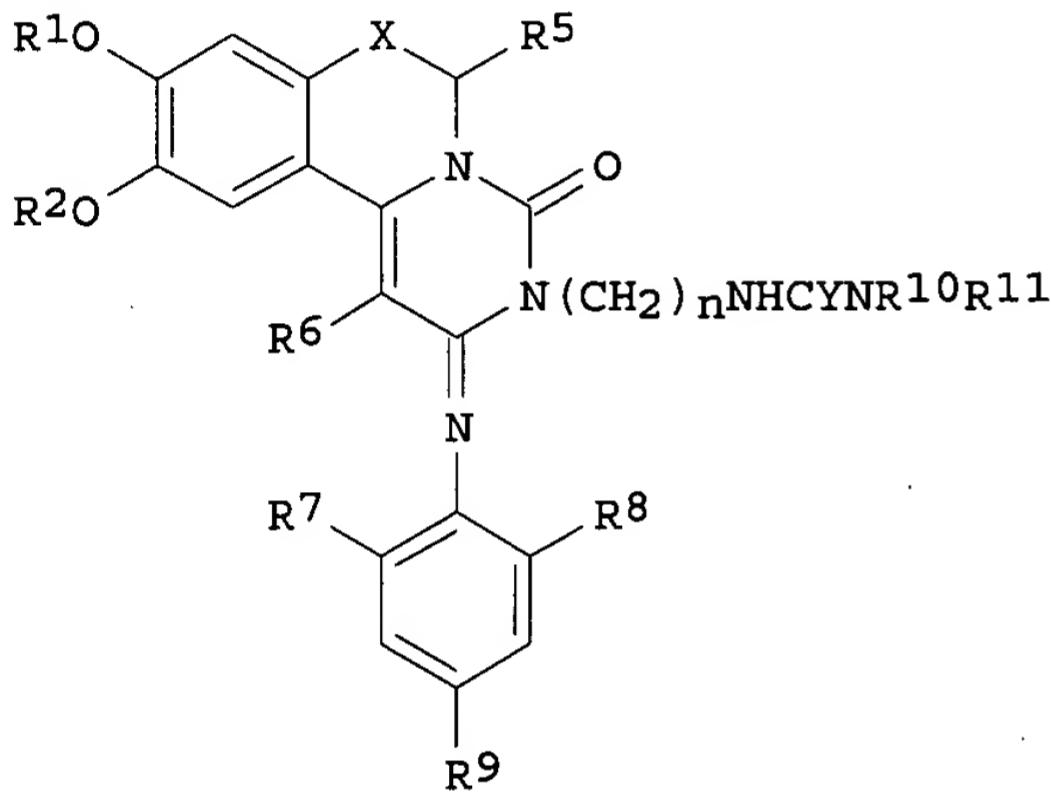
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WO 2000058308	A1	20001005	WO 2000-GB1193	20000329
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RW:	GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE,			

DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF,
 CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG
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 AU 2000041274 A5 20001016 AU 2000-41274 20000329
 EP 1165558 A1 20020102 EP 2000-920857 20000329
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 IE, SI, LT, LV, FI, RO
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 JP 2002540207 T2 20021126 JP 2000-608010 20000329
 US 2003036542 A1 20030220 US 2001-964260 20010926
 NO 2001004728 A 20011123 NO 2001-4728 20010928
 PRIORITY APPLN. INFO.: GB 1999-7454 A 19990331
 GB 1999-9802 A 19990428
 WO 2000-GB1193 W 20000329

OTHER SOURCE(S) :

GI

OTHER SOURCE(S) : MARPAT 133:266869



AB Title compds. [I; R₁, R₂ = alkyl, acyl; R₅ = H, alkyl, alkenyl, alkynyl; R₆ = H, alkyl, alkenyl, alkynyl, amino, alkylamino, dialkylamino, acylamino; R₇, R₈ = H, halo, OH, CF₃, alkyl, alkenyl, alkynyl, acyl, alkythio, alkoxy, cycloalkyl; R₉ = H, halo, OH, CF₃, alkyl, alkenyl, alkynyl, acyl, alkythio, alkoxy, cycloalkyl; X = OCH₂, CR₃R₄; R₃, R₄ = H, alkyl; R₁₀, R₁₁ = H, alkyl, cycloalkyl, Ph; Y = O, CHNO₂, NCN, NH, NNO₂; n = 2-4], were prepd. I have a longer duration of action than the known compd. trequinsin (9,10-dimethoxy-3-methyl-2-mesitylimino-2,3,6,7-tetrahydro-4H-pyrimido[6,1-a]isoquinolin-4-one) and do not have trequinsin's very bitter taste. Thus, Na cyanate was added dropwise to 9,10-dimethoxy-2-(2,4,6-trimethylphenylimino)-3-(2-aminoethyl)-3,4,6,7-tetrahydro-2H-pyrimido[6,1-a]isoquinolin-4-one (prepn. given) in aq. HCl at 80.degree. followed by stirring for 2 h to give 54% 9,10-dimethoxy-2-(2,4,6-trimethylphenylimino)-3-(N-carbamoyl-2-aminoethyl)-3,4,6,7-tetrahydro-2H-pyrimido[6,1-a]isoquinolin-4-one (II). II inhibited PDE3 with IC₅₀ = 0.46 .mu.M and was tasteless.

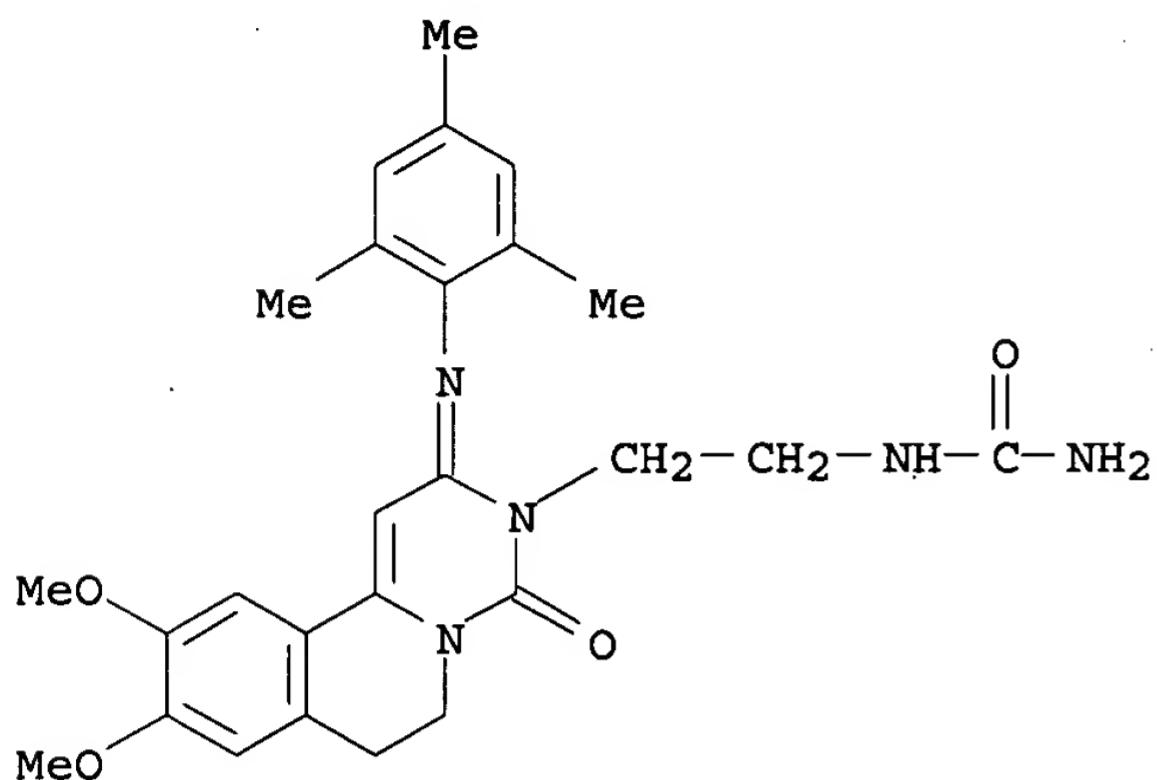
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298680-31-6P 298680-32-7P 298680-33-8P
298680-34-9P 298680-35-0P 298680-36-1P
298680-37-2P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of 2-aryliminopyrimido[6,1-a]isoquinolin-4-ones as phosphodiesterase inhibitors)

RN 298680-25-8 CAPLUS

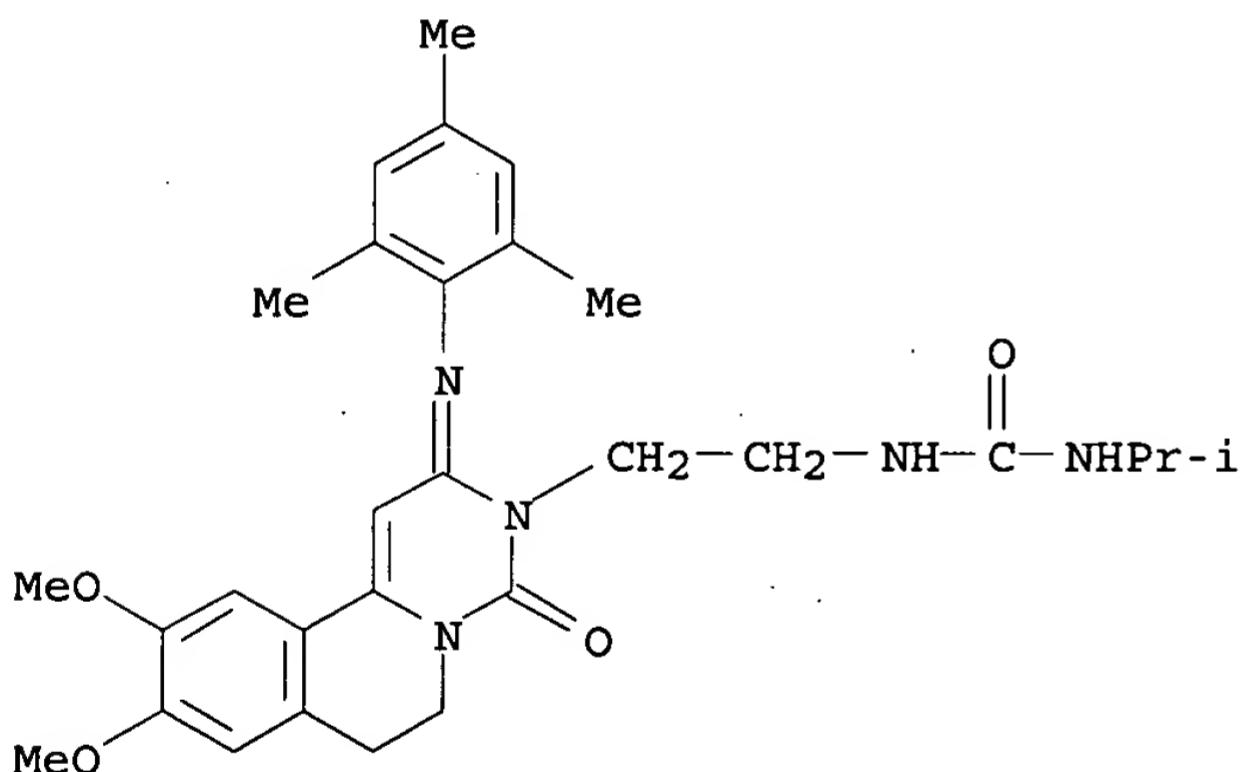
09/ 964,260

CN Urea, [2- [6,7-dihydro-9,10-dimethoxy-4-oxo-2- [(2,4,6-trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl] - (9CI) (CA INDEX NAME)



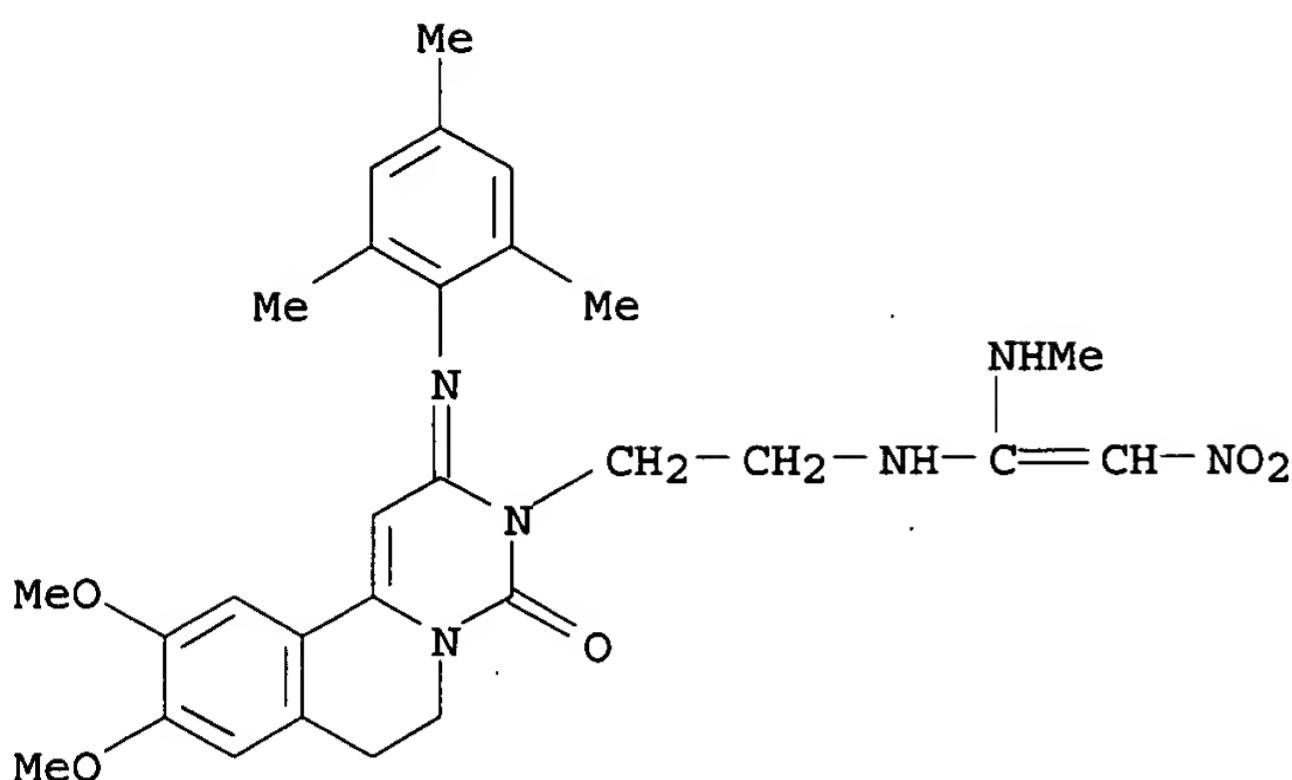
RN 298680-26-9 CAPLUS

CN Urea, N- [2- [6,7-dihydro-9,10-dimethoxy-4-oxo-2- [(2,4,6-trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl] -N' - (1-methylethyl) - (9CI) (CA INDEX NAME)

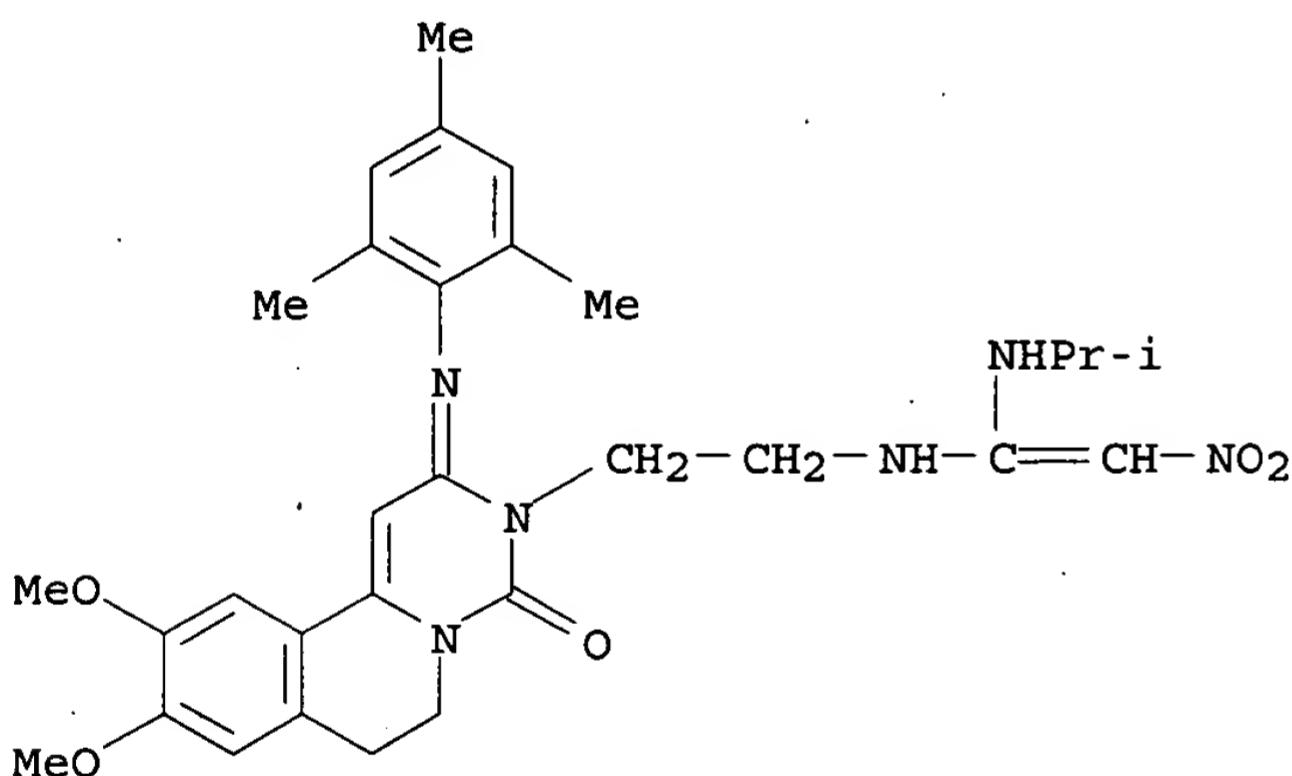


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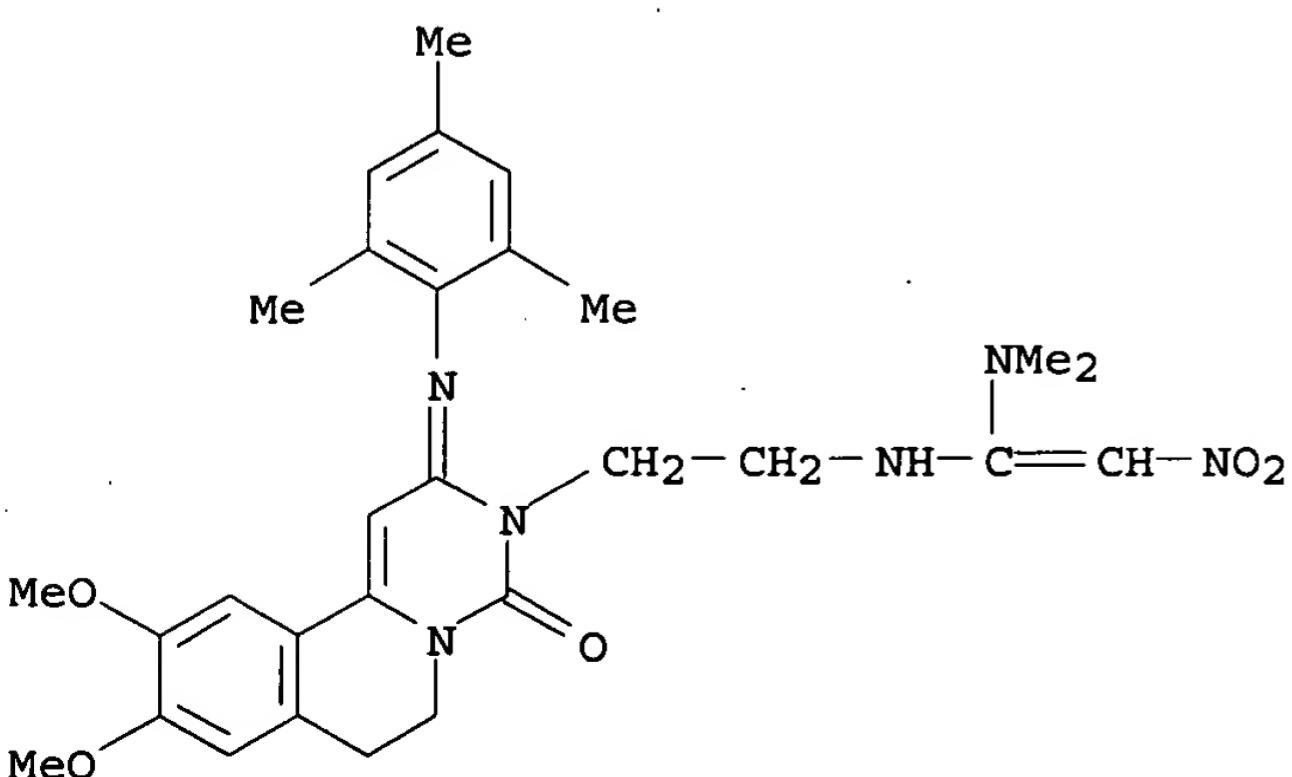
CN 4H-Pyrimido[6,1-a]isoquinolin-4-one, 2,3,6,7-tetrahydro-9,10-dimethoxy-3-[2-[(1-methylamino)-2-nitroethenyl]amino]ethyl]-2-[(2,4,6-trimethylphenyl)imino] - (9CI) (CA INDEX NAME)



RN 298680-28-1 CAPLUS
CN 4H-Pyrimido[6,1-a]isoquinolin-4-one, 2,3,6,7-tetrahydro-9,10-dimethoxy-3-[2-[(1-methylethyl)amino]-2-nitroethyl]amino]ethyl]-2-[(2,4,6-trimethylphenyl)imino]- (9CI) (CA INDEX NAME)

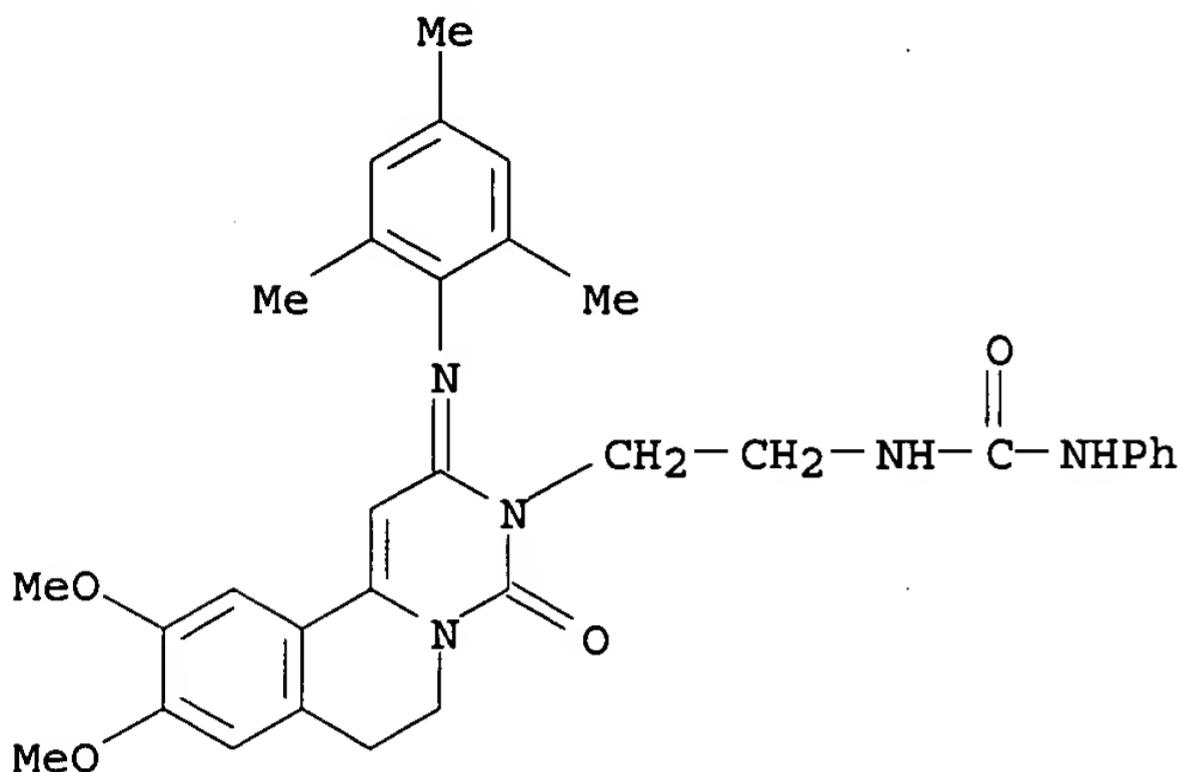


RN 298680-29-2 CAPLUS
CN 4H-Pyrimido[6,1-a]isoquinolin-4-one, 3-[2-[(dimethylamino)-2-nitroethyl]amino]ethyl]-2,3,6,7-tetrahydro-9,10-dimethoxy-2-[(2,4,6-trimethylphenyl)imino]- (9CI) (CA INDEX NAME)



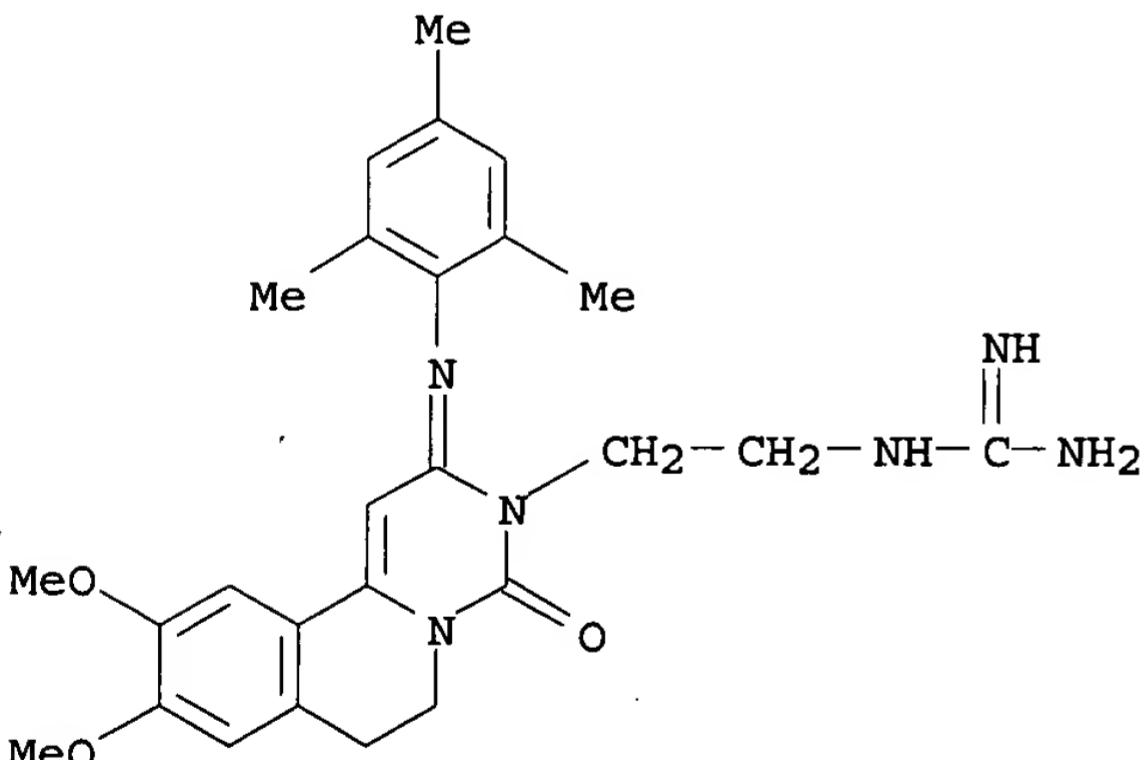
RN 298680-30-5 CAPLUS

CN Urea, N-[2-[6,7-dihydro-9,10-dimethoxy-4-oxo-2-[(2,4,6-trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl]-N'-phenyl- (9CI) (CA INDEX NAME)



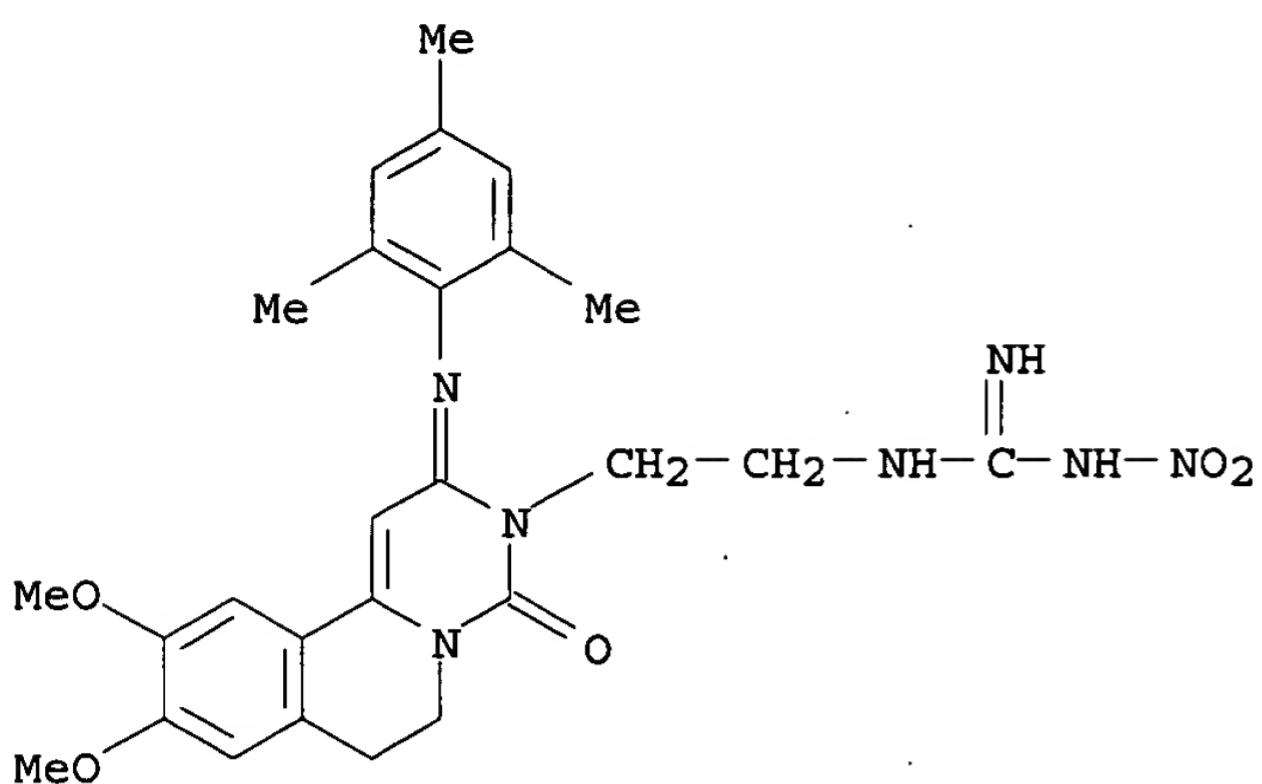
RN 298680-31-6 CAPLUS

CN Guanidine, [2-[6,7-dihydro-9,10-dimethoxy-4-oxo-2-[(2,4,6-trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl]- (9CI) (CA INDEX NAME)

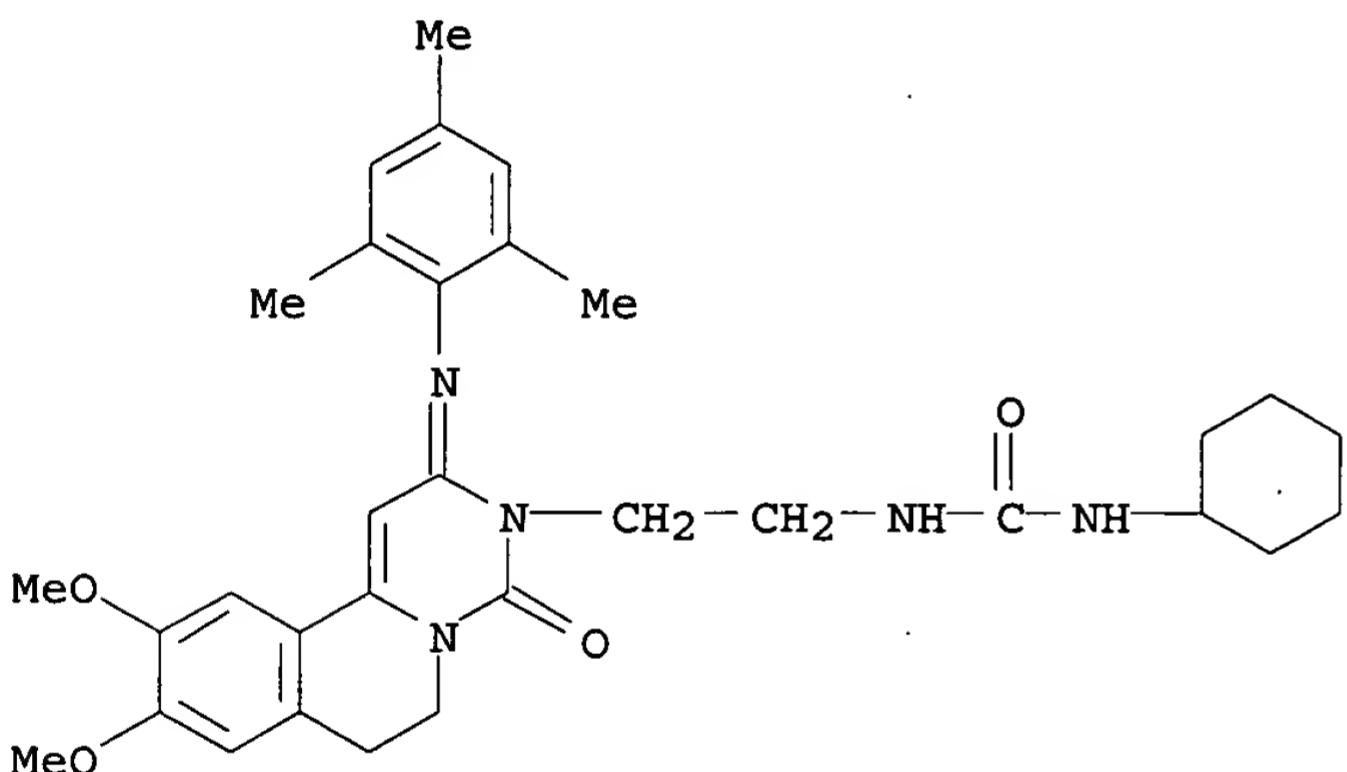


RN 298680-32-7 CAPLUS

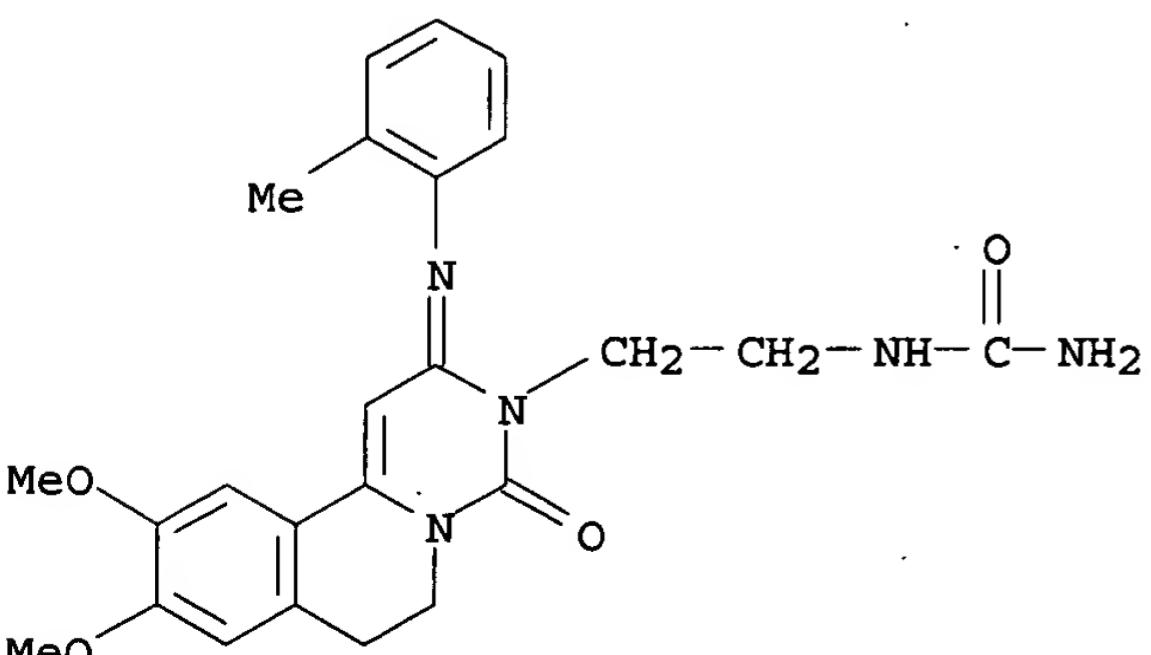
CN Guanidine, N-[2-[6,7-dihydro-9,10-dimethoxy-4-oxo-2-[(2,4,6-trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl]-N'-nitro- (9CI) (CA INDEX NAME)



RN 298680-33-8 CAPLUS
CN Urea, N-cyclohexyl-N'-[2-[6,7-dihydro-9,10-dimethoxy-4-oxo-2-[(2,4,6-trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl] - (9CI) (CA INDEX NAME)

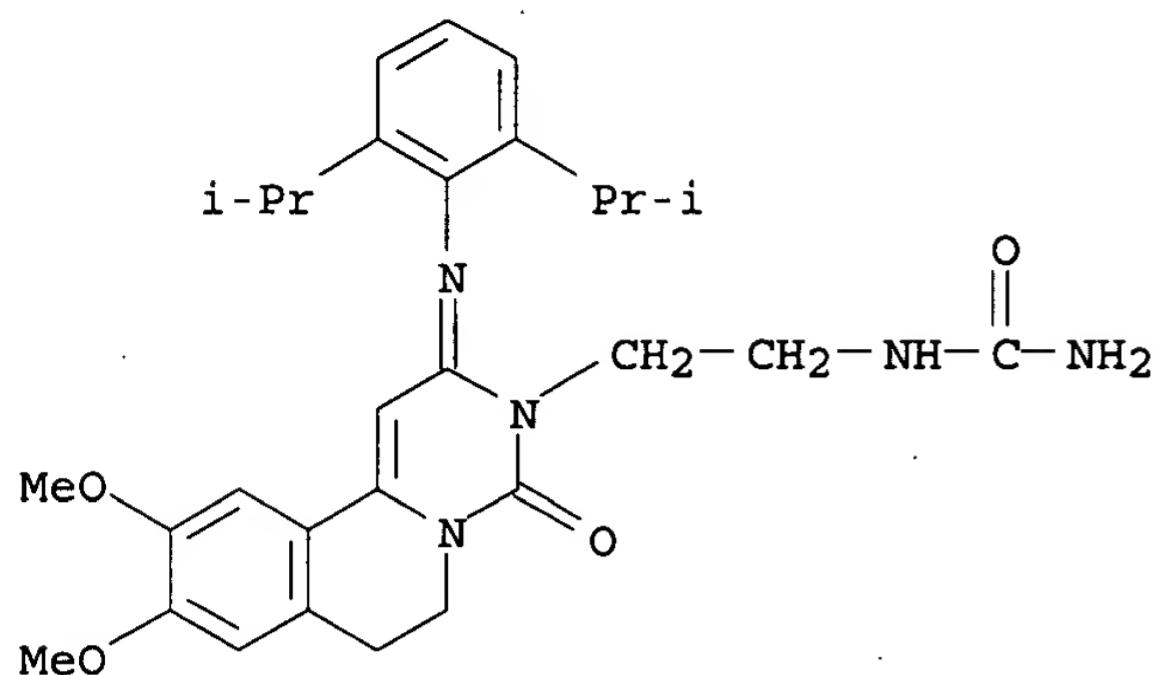


RN 298680-34-9 CAPLUS
CN Urea, [2-[6,7-dihydro-9,10-dimethoxy-2-[(2-methylphenyl)imino]-4-oxo-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl] - (9CI) (CA INDEX NAME)



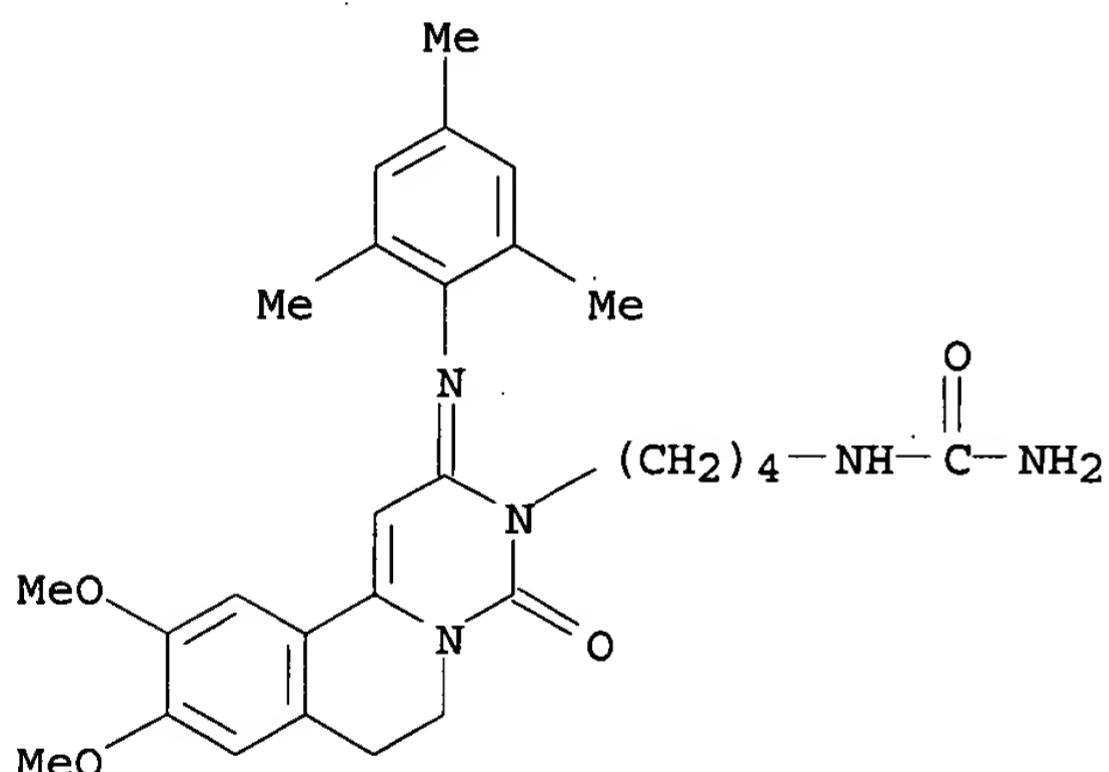
RN 298680-35-0 CAPLUS
CN Urea, [2-[2-[[2,6-bis(1-methylethyl)phenyl]imino]-6,7-dihydro-9,10-

dimethoxy-4-oxo-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl] - (9CI) (CA INDEX NAME)



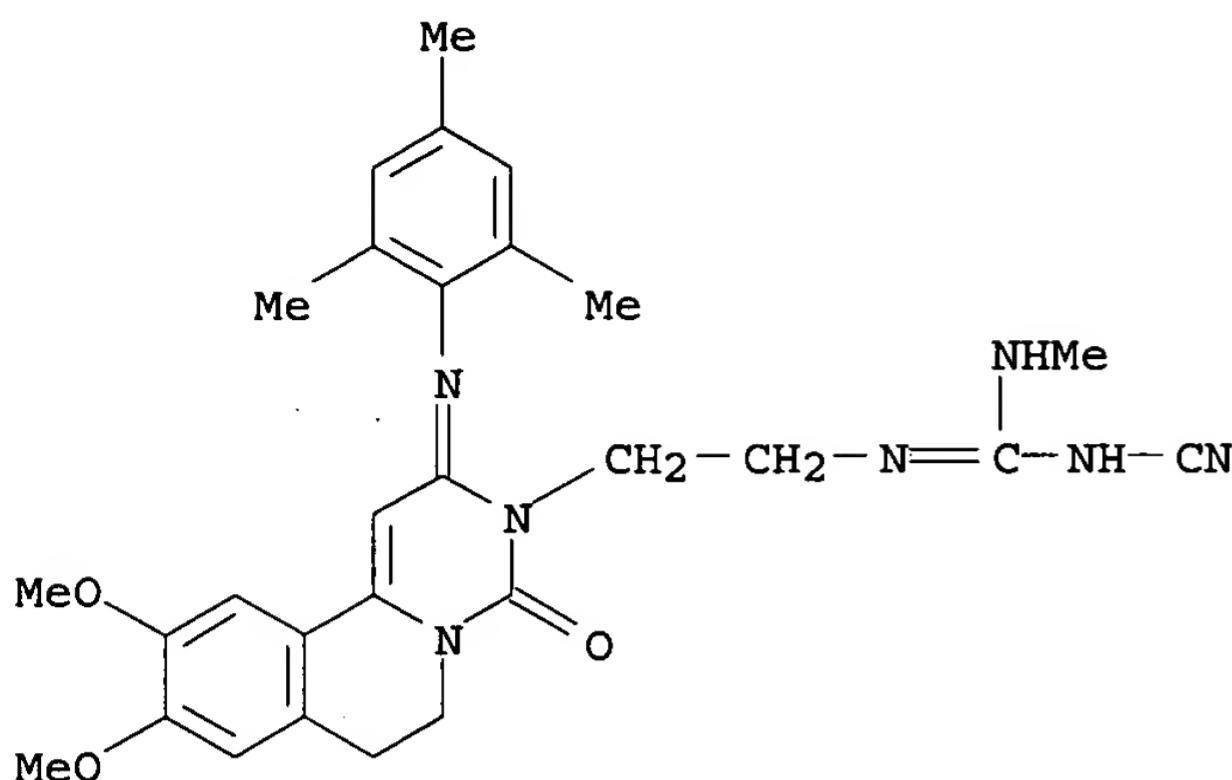
RN 298680-36-1 CAPLUS

CN Urea, [4- [6,7-dihydro-9,10-dimethoxy-4-oxo-2- [(2,4,6-trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]butyl] - (9CI) (CA INDEX NAME)



RN 298680-37-2 CAPLUS

CN Guanidine, N-cyano-N' - [2- [6,7-dihydro-9,10-dimethoxy-4-oxo-2- [(2,4,6-trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl] -N' - methyl - (9CI) (CA INDEX NAME)



IT 298680-40-7P

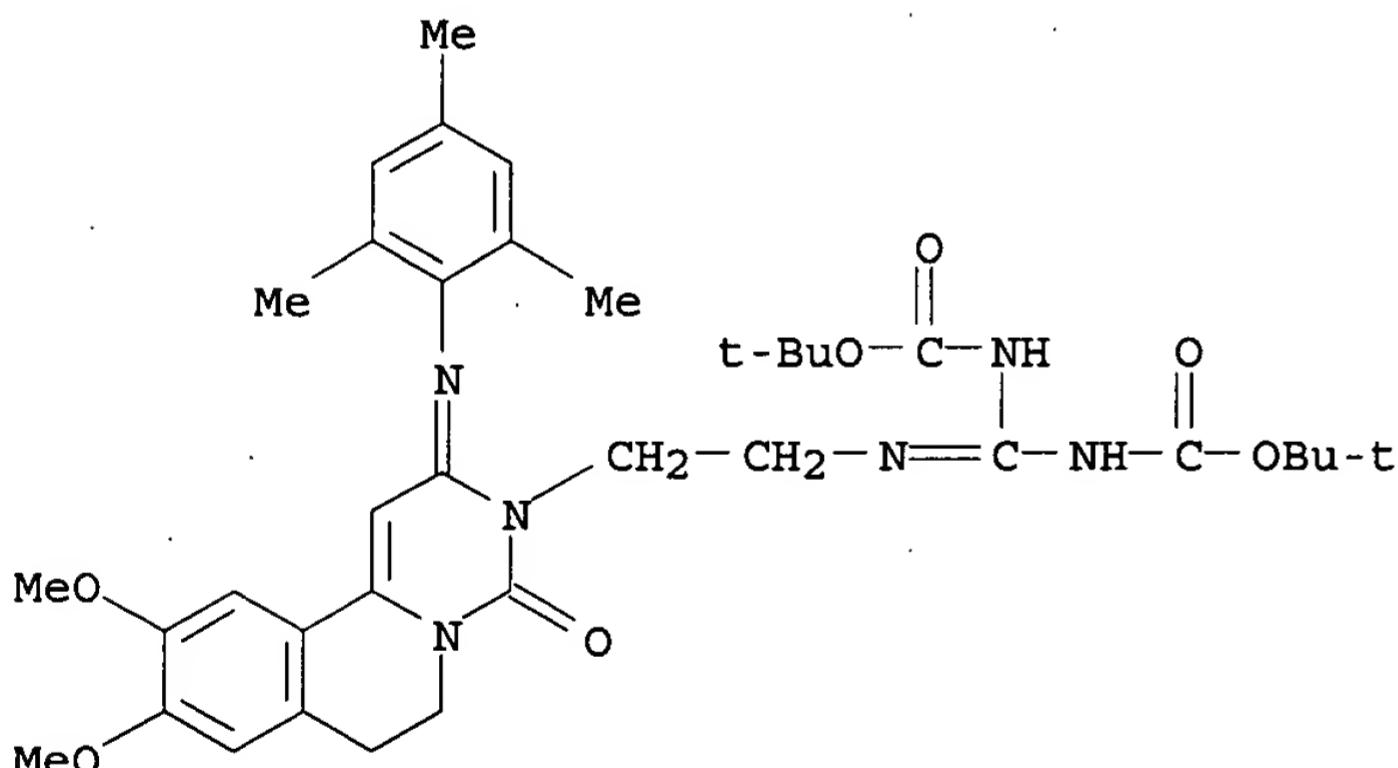
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(prepn. of 2-aryliminopyrimido[6,1-a]isoquinolin-4-ones as phosphodiesterase inhibitors)

RN 298680-40-7 CAPLUS

CN Carbamic acid, [[2-[6,7-dihydro-9,10-dimethoxy-4-oxo-2-[(2,4,6-

trimethylphenyl)imino]-2H-pyrimido[6,1-a]isoquinolin-3(4H)-yl]ethyl]carbonimidoyl]bis-, bis(1,1-dimethylethyl) ester (9CI) (CA INDEX NAME)



REFERENCE COUNT:

1

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SINCE FILE
ENTRY

TOTAL SESSION

DISCOUNT AMOUNTS (F)

ENTRY

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FILE COVERS 1907-1966

FILE LAST UPDATED: 01 May 1997 (19970501/UP)

This file contains CAS Registry Numbers for easy and accurate substance identification. Title keywords, authors, patent assignees, and patent information, e.g., patent numbers, are now searchable from 1907-1966. TIFF images of CA abstracts printed between 1907-1966 are available in the PAGE display formats.

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FILE 'CAPLUS' ENTERED AT 14:20:36 ON 15 AUG 2003

L3 1 S L2

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L4 0 L2

=> d his

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FILE 'REGISTRY' ENTERED AT 14:20:04 ON 15 AUG 2003

L1 STRUCTURE uploaded
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FILE 'CAPLUS' ENTERED AT 14:20:36 ON 15 AUG 2003

L3 1 S L2

FILE 'CAOLD' ENTERED AT 14:21:07 ON 15 AUG 2003

L4 0 S L2

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